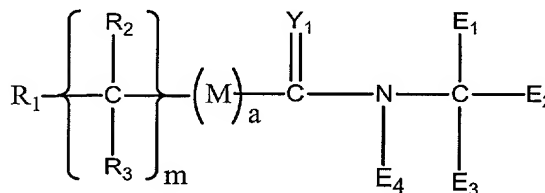


WHAT IS CLAIMED IS:

1. A compound comprising the formula:

(I)



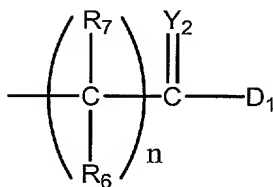
wherein:

R_1 is a polymeric residue;

Y_1 is O, S or NR_4 ;

M is O, S or NR_5 ;

E_1 is



$E_{2,4}$ are independently H, E_1 or

(a) is zero or one;

(m) is zero or a positive integer;

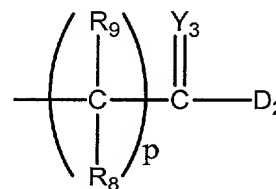
(n) and (p) are independently 0 or a positive integer;

$Y_{2,3}$ are independently O, S or NR_{10} ;

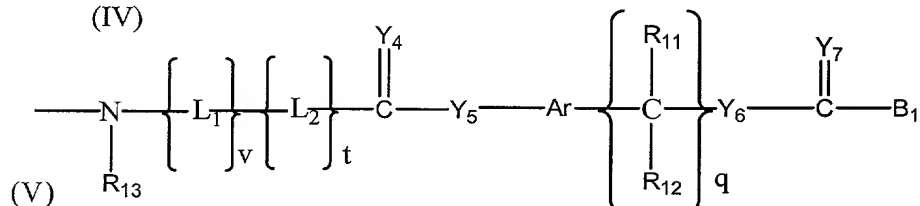
R_{2-10} are independently selected from the group consisting of hydrogen,

C_{1-6} alkyls, C_{3-12} branched alkyls, C_{3-8} cycloalkyls, C_{1-6} substituted alkyls, C_{3-8} substituted cycloalkyls, aryls, substituted aryls, aralkyls, C_{1-6} heteroalkyls, substituted C_{1-6} heteroalkyls, C_{1-6} alkoxy, phenoxy and C_{1-6} heteroalkoxy;

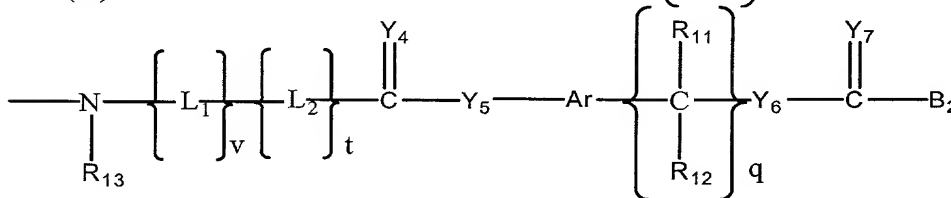
D_1 and D_2 are independently OH,



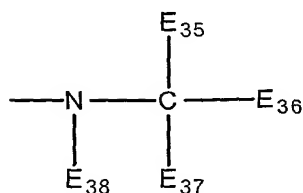
(IV)



(V)

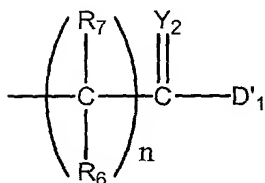


4. The compound of claim 1, wherein said terminal branching group comprises the formula:

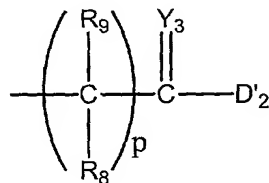


wherein

E₃₅ is



E₃₆₋₃₈ are independently H, E₃₅ or

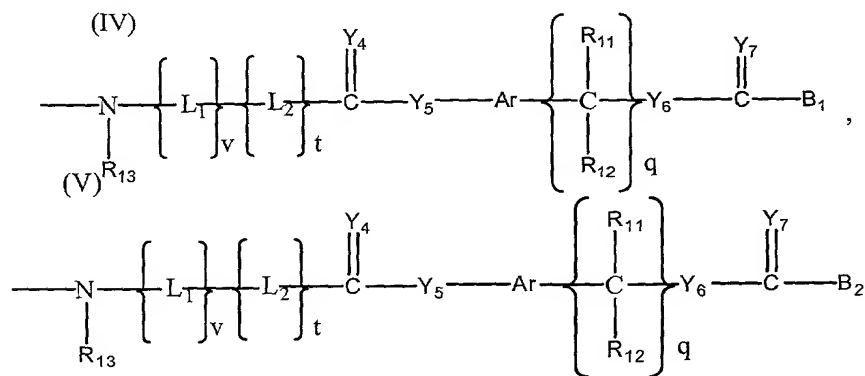


(n) and (p) are independently 0 or a positive integer;

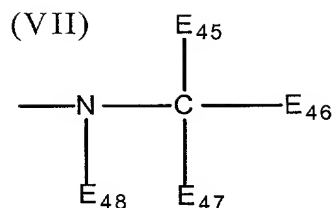
Y_{2,3} are independently O, S or NR₁₀;

R₆₋₁₀ are independently selected from the group consisting of hydrogen, C₁₋₆ alkyls, C₃₋₁₂ branched alkyls, C₃₋₈ cycloalkyls, C₁₋₆ substituted alkyls, C₃₋₈ substituted cycloalkyls, aryls, substituted aryls, aralkyls, C₁₋₆ heteroalkyls, substituted C₁₋₆ heteroalkyls, C₁₋₆ alkoxy, phenoxy and C₁₋₆ heteroalkoxy;

D'₁ and D'₂ are independently OH,



or



wherein (v) and (t) are independently 0 or a positive integer up to about 6;

(q) is zero or a positive integer;

L_1 and L_2 are independently selected bifunctional linkers;

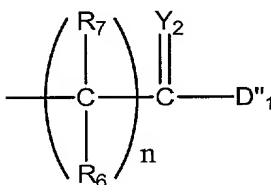
Y_{4-7} are independently selected from the group consisting of O, S and NR_{14} ;

R_{11-14} are independently selected from the group consisting of hydrogen, C_{1-6} alkyls, C_{3-12} branched alkyls, C_{3-8} cycloalkyls, C_{1-6} substituted alkyls, C_{3-8} substituted cycloalkyls, aryls, substituted aryls, aralkyls, C_{1-6} heteroalkyls, substituted C_{1-6} heteroalkyls, C_{1-6} alkoxy, phenoxy and C_{1-6} heteroalkoxy;

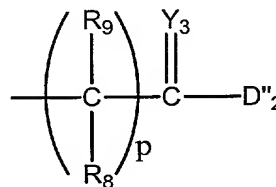
Ar is a moiety which when included in Formula (I) forms a multi-substituted aromatic hydrocarbon or a multi-substituted heterocyclic group;

B_1 and B_2 are independently selected from the group consisting of leaving groups, OH, residues of hydroxyl-containing moieties or amine-containing moieties;

E_{45} is

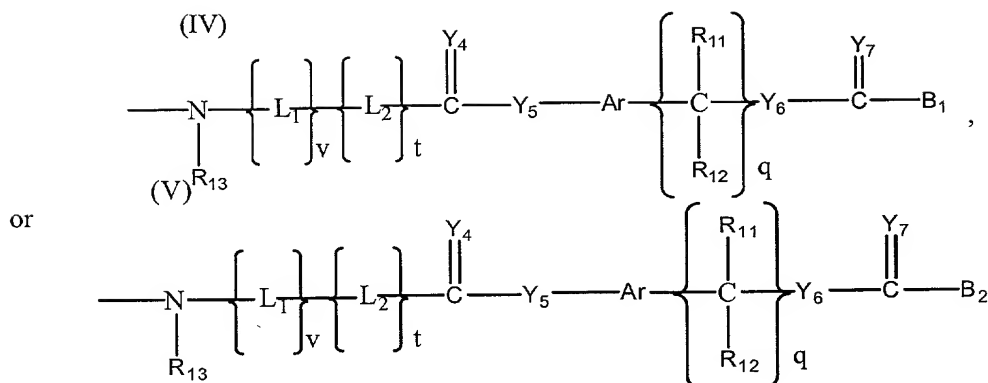


E_{46-48} are independently H, E_{45} or



wherein

D''_1 and D''_2 are independently OH,



5. The compound of claim 3, Y₁ is O.
6. The compound of claim 1, wherein R₁ comprises a polyalkylene oxide residue.
7. The compound of claim 6, wherein R₁ comprises a polyethylene glycol residue.
8. The compound of claim 3, wherein R₁ comprises a polyethylene glycol residue.
9. The compound of claim 6, wherein R₁ is selected from the group consisting of

-C(=Y₈)-(CH₂)_f-O-(CH₂CH₂O)_x-A, -C(=Y₈)-Y₉-(CH₂)_f-O-(CH₂CH₂O)_x-A,
 -C(=Y₈)-NR₂₀-(CH₂)_f-O-(CH₂CH₂O)_x-A, -(CR₂₁R₂₂)_e-O-(CH₂)_f-O-(CH₂CH₂O)_x-A,
 -NR₂₀-(CH₂)_f-O-(CH₂CH₂O)_x-A, -C(=Y₈)-(CH₂)_f-O-(CH₂CH₂O)_x-(CH₂)_f-C(=Y₈)-,
 -C(=Y₈)-Y₉-(CH₂)_f-O-(CH₂CH₂O)_x-(CH₂)_f-Y₉-C(=Y₈)-,
 -C(=Y₈)-NR₂₀-(CH₂)_f-O-(CH₂CH₂O)_x-(CH₂)_f-NR₂₀-C(=Y₈)-,
 -(CR₂₁R₂₂)_e-O-(CH₂)_f-O-(CH₂CH₂O)_x-(CH₂)_f-O-(CR₂₁R₂₂)_e-, and
 -NR₂₀-(CH₂)_f-O-(CH₂CH₂O)_x-(CH₂)_f-NR₂₀-

wherein:

Y₈ and Y₉ are independently O, S or NR₂₀;

x is the degree of polymerization;

R₂₀, R₂₁ and R₂₂ are independently selected from among H, C₁₋₆ alkyls,

C₃₋₁₂ branched alkyls, C₃₋₈ cycloalkyls, C₁₋₆ substituted alkyls, C₃₋₈ substituted cycloalkyls,
 aryls, substituted aryls, aralkyls, C₁₋₆ heteroalkyls, substituted C₁₋₆ heteroalkyls,

C₁₋₆ alkoxy, phenoxy and C₁₋₆ heteroalkoxy;

e and f are independently zero, one or two; and

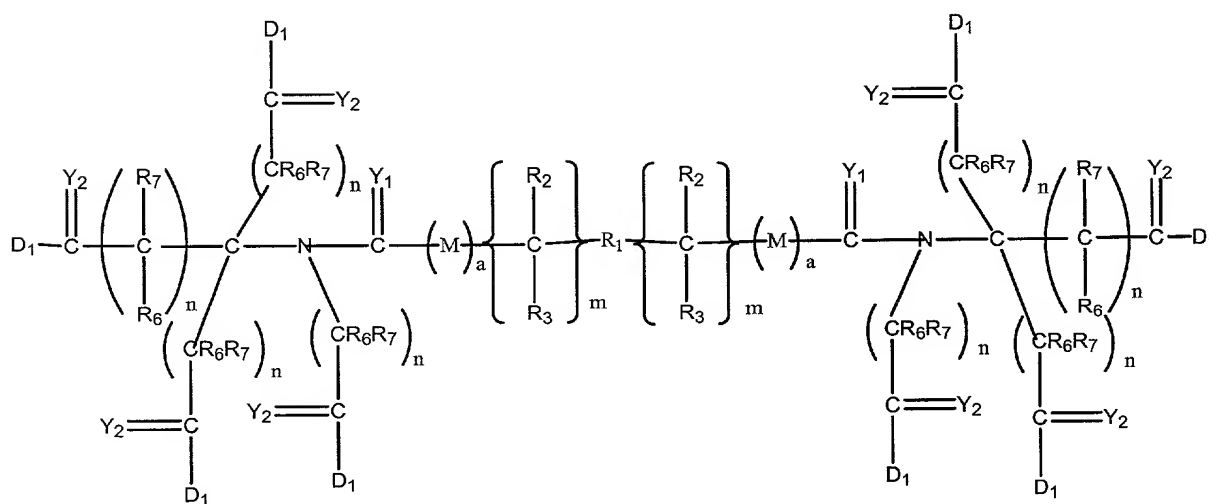
A is a capping group.

10. The compound of claim 9, wherein R_1 comprises $-O-(CH_2CH_2O)_x$ and x is a positive integer so that the weight average molecular weight is at least about 20,000.

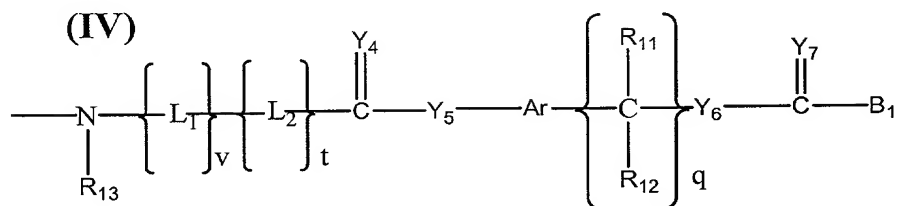
11. The compound of claim 3, wherein R_1 has a weight average molecular weight of from about 20,000 to about 100,000.

12. The compound of claim 3, wherein R_1 has a weight average molecular weight of from about 25,000 to about 60,000.

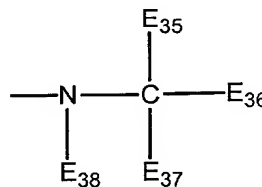
13. A compound of claim 3, comprising the formula



14. The compound of claim 13, wherein D_1 is



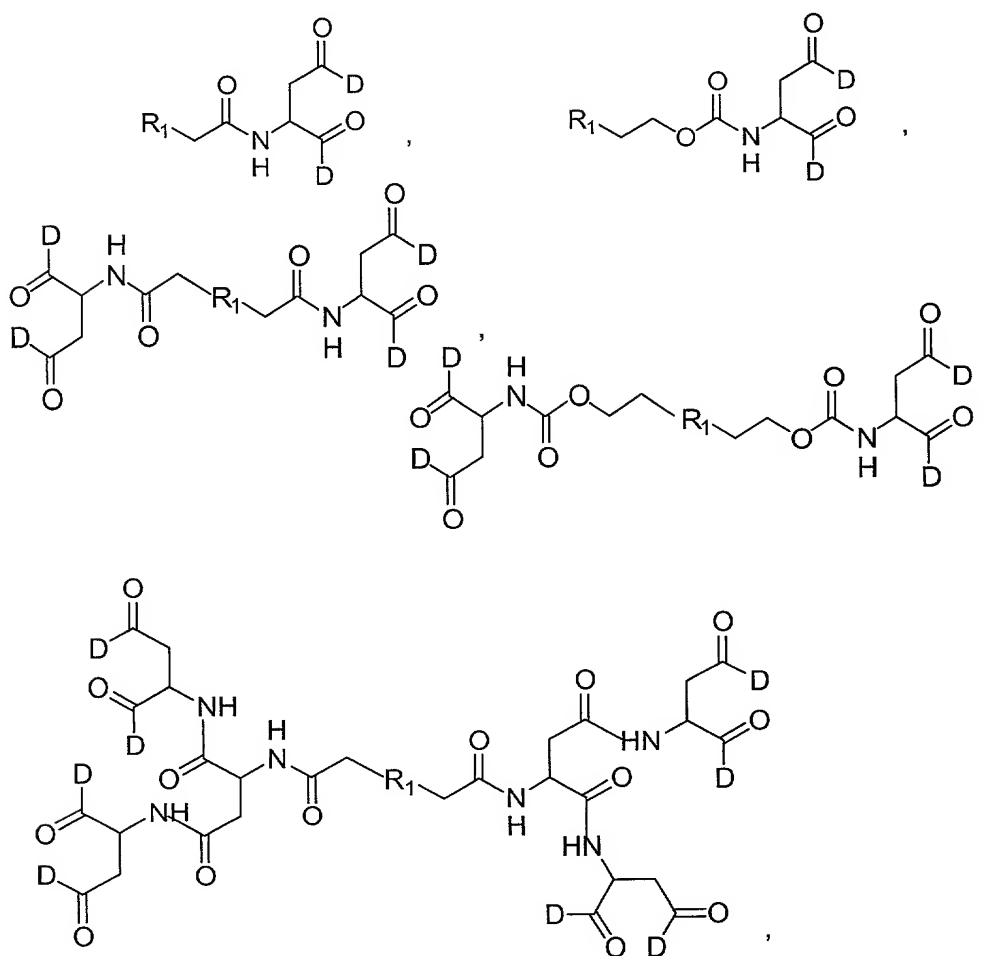
15. The compound of claim 13, wherein D_1 is

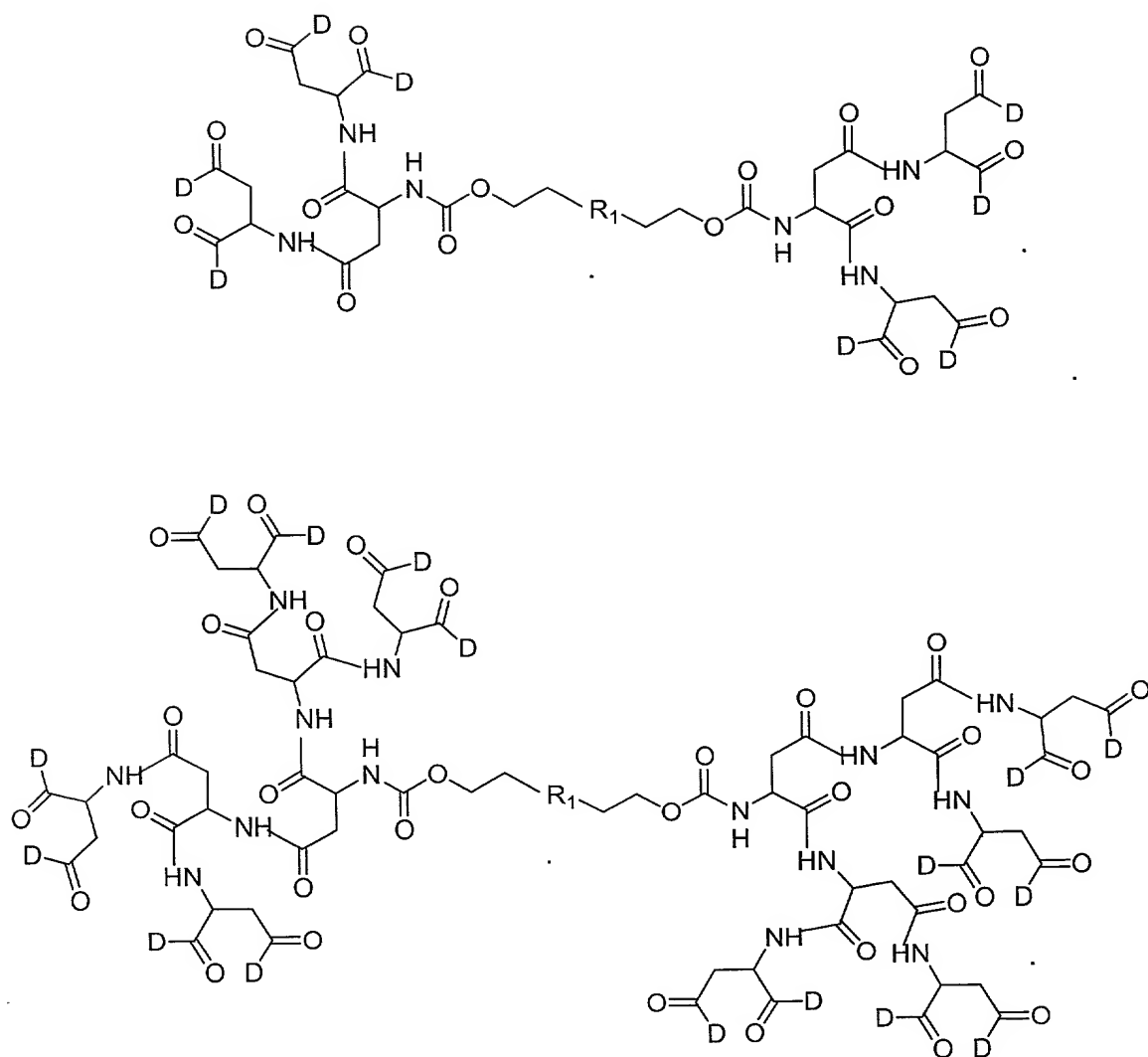


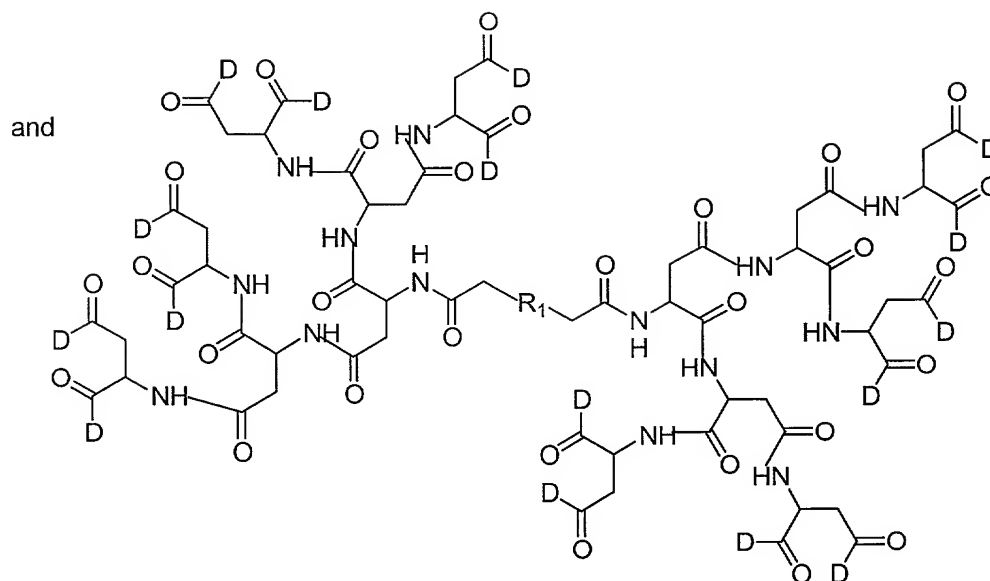
16. The compound of claim 1, wherein L_1 is $(CH_2CH_2O)_2$.

17. The compound of claim 1, wherein L_2 is selected from the group consisting of $-CH_2-$, $-CH(CH_3)-$, $-CH_2C(O)NHCH(CH_3)-$, $-(CH_2)_2-$, $-CH_2C(O)NHCH_2-$, $-(CH_2)_2NH-$, $-(CH_2)_2NH-C(O)(CH_2)_2NH-$ and $-CH_2C(O)NHCH(CH_2CH(CH_3)_2)-$.

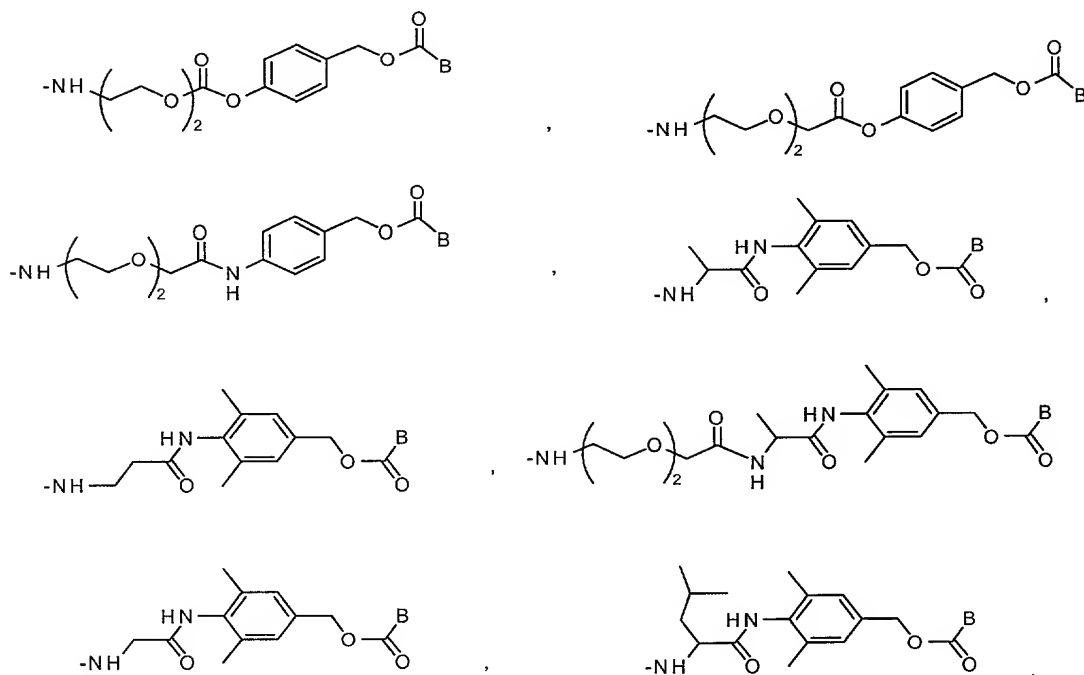
18. A compound of claim 1, selected from the group consisting of:

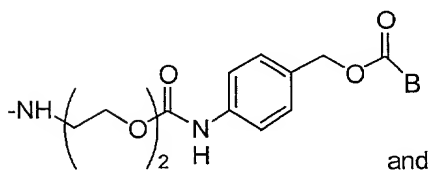
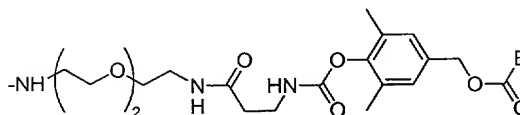
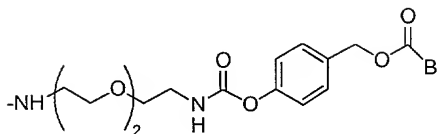
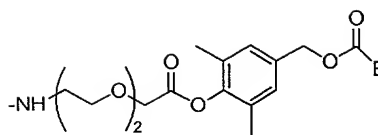
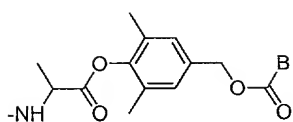




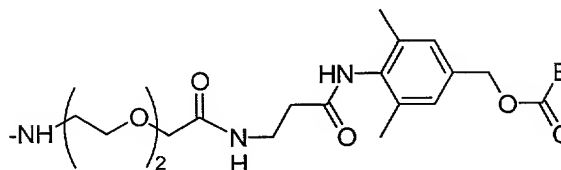


wherein R_1 is a PEG residue and D is selected from the group comprising:





and



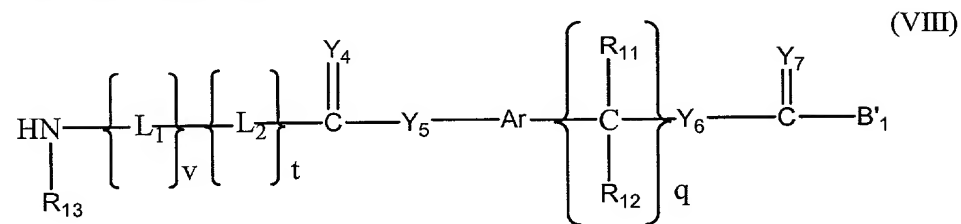
where B is a residue of an amine or a hydroxyl- containing drug.

19. A compound of claim 18, wherein B is a residue of a member of the group consisting of: daunorubicin, doxorubicin; *p*-aminoaniline mustard, melphalan, Ara-C (cytosine arabinoside), leucine-Ara-C, and gemcitabine

20. A method of treatment, comprising administering to a mammal in need of such treatment an effective amount of a compound of claim 1, wherein D₁ is a residue of a biologically active moiety.

21. A method of treatment, comprising administering to a mammal in need of such treatment an effective amount of a compound of claim 18.

22. A method of preparing a polymer conjugate, comprising:
reacting a compound of the formula (VIII):



wherein (v) and (t) are independently 0 or a positive integer up to about 6;

L_1 and L_2 are independently selected bifunctional linkers;

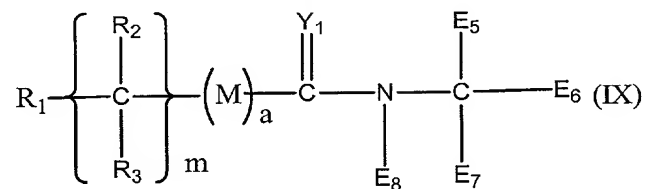
Y_{4-7} are independently selected from the group consisting of O, S and NR_{14} ;

R_{11-14} are independently selected from the group consisting of hydrogen, C_{1-6} alkyls, C_{3-12} branched alkyls, C_{3-8} cycloalkyls, C_{1-6} substituted alkyls, C_{3-8} substituted cycloalkyls, aryls, substituted aryls, aralkyls, C_{1-6} heteroalkyls, substituted C_{1-6} heteroalkyls, C_{1-6} alkoxy, phenoxy and C_{1-6} heteroalkoxy;

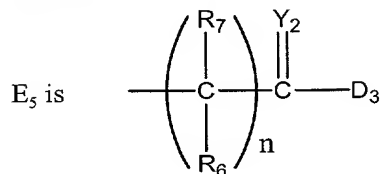
Ar is a moiety which when included in Formula (I) forms a multi-substituted aromatic hydrocarbon or a multi-substituted heterocyclic group; and

B'_1 is a residue of a hydroxyl- or an amine-containing moiety;

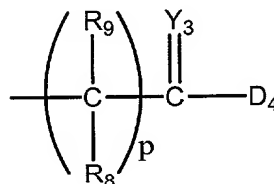
with a compound of the formula (IX):



wherein



E_{6-8} are independently H, E_5 or



wherein

D_3 and D_4 are independently OH, a leaving group which is capable of reacting with an unprotected amine or hydroxyl or a terminal branching group;

R_1 is a polymeric residue;

Y_1 is O, S or NR_4 ;

M is O, S or NR_5 ;

(n) and (p) are independently 0 or a positive integer;

Y_{2-3} are independently O, S or NR_{10} ; and

R_{2-10} are independently selected from the group consisting of hydrogen, C_{1-6} alkyls, C_{3-12} branched alkyls, C_{3-8} cycloalkyls, C_{1-6} substituted alkyls, C_{3-8} substituted cycloalkyls, aryls, substituted aryls, aralkyls, C_{1-6} heteroalkyls, substituted C_{1-6} heteroalkyls, C_{1-6} alkoxy, phenoxy and C_{1-6} heteroalkoxy;

under conditions sufficient to cause a polymeric conjugate to be formed.